

REMARKS

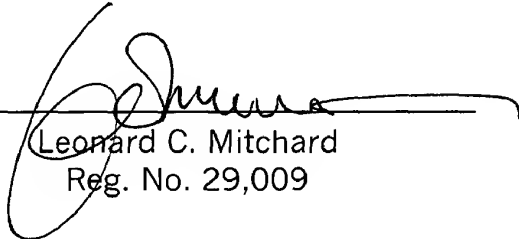
The above amendments have been made to place the application in a more traditional format.

Attached hereto is a marked-up version of the changes made to the specification and claims by the current amendment. The attached pages are captioned "**Version With Markings To Show Changes Made.**"

Respectfully submitted,

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By: _____


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VERSION WITH MARKINGS TO SHOW CHANGES MADE

IN THE CLAIMS

5. (Amended) A compound according to claim 2 [or claim 4] wherein R^1 , R^2 and R^3 are each methyl.
6. (Amended) A compound according to [any one of claims 2, 4 or 5] claim 2 wherein R^4 is hydrogen.
7. (Amended) A compound according to [any one of claims 2, 4, 5 or 6] claim 2 wherein R^6 is hydrogen, halogeno, amino, carboxy, hydroxy, C_{1-7} alkoxy or a group Y^4R^{35} (wherein Y^4 is $-C(O)-$, $-O-$ or $-OSO_2-$ and R^{35} is C_{1-7} alkyl, C_{1-7} alkoxy (which alkyl or alkoxy may bear one or more substituents selected from halogeno), R^{48} (wherein R^{48} is a benzyl group) or R^{53} (wherein R^{53} is a 5-6-membered saturated heterocyclic group (linked via carbon or nitrogen) with 1-2 heteroatoms selected independently from O, S and N)).
8. (Amended) A compound according to [any one of claims 2, 4, 5, 6 or 7] claim 2 wherein R^6 is hydrogen, $C(O)OCH_3$ or methoxy.
9. (Amended) A compound according to [any one of claims 2, 4, 5, 6, 7 or 8] claim 2 wherein R^5 is hydrogen, halogeno, amino, carboxy, carbamoyl,

C₁₋₇alkanoyl, C₁₋₇thioalkoxy, or a group -Y⁴R³⁵

(wherein Y⁴ is -C(O)-, -OC(O)-, -O-, -SO-, -OSO₂-, -NR³⁶-, -NR³⁷C(O)- or -C(O)NR³⁸.

(wherein R³⁶, R³⁷ and R³⁸, which may be the same or different, each represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₁₋₃alkyl) and

R³⁵ is a sugar moiety, a mono-peptide, a di-peptide, a tri-peptide, a tetra-peptide, C₁₋₇alkyl, C₁₋₇alkoxy, C₁₋₇alkanoyl, C₁₋₇alkanoylaminoC₁₋₇alkyl,

(which alkyl, alkoxy, alkanoyl, alkanoylaminoalkyl may bear one or more substituents selected from:

halogeno, amino, hydroxy, carboxy, and a group -Y⁵R⁴⁰ (wherein Y⁵ is -C(O)-O- or -O-C(O)- and R⁴⁰ is C₁₋₇alkyl or a group R⁴³ wherein R⁴³ is a benzyl group),

R⁴⁸ (wherein R⁴⁸ is a phenyl group, a benzyl group or a 5-10-membered aromatic heterocyclic group (linked via carbon or nitrogen) with 1-4 heteroatoms selected independently from O, N and S, which phenyl, benzyl or aromatic heterocyclic group may bear one or more substituents selected from

hydroxy, fluoro, amino, C₁₋₄alkoxy, C₁₋₄hydroxyalkyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, di(C₁₋₄hydroxyalkyl)amino C₁₋₄alkyl, di(C₁₋₄aminoalkylamino C₁₋₄alkyl, C₁₋₄hydroxyalkoxy, carboxy, C₁₋₄carboxyalkyl, cyano, -CONR⁴⁹R⁵⁰, -NR⁵¹COR⁵² (wherein R⁴⁹, R⁵⁰, R⁵¹ and R⁵², which may be the same or different, each represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and C₁₋₄alkylR⁵³ (wherein R⁵³ is as defined herein),

$C_{1-7}alkylR^{48}$ (wherein R^{48} is as defined herein),

R^{53} (wherein R^{53} is a 5-6-membered saturated heterocyclic group (linked via carbon or nitrogen) with 1-2 heteroatoms, selected independently from O, S and N, which heterocyclic group may bear 1 or 2 substituents selected from

oxo, hydroxy, fluoro, chloro, alkyl, $C_{1-4}hydroxyalkyl$, $C_{1-4}alkoxy$,

$C_{1-4}carboxyalkyl$, $C_{1-4}aminoalkyl$, $di(C_{1-4}alkyl)amino$ $C_{1-4}alkyl$, $C_{1-4}alkoxy$ $C_{1-4}alkyl$,

$C_{1-4}alkylsulphonyl$ $C_{1-4}alkyl$ and R^{54} (wherein R^{54} is a 5-6-membered saturated heterocyclic group (linked via carbon or nitrogen) with 1-2 heteroatoms, selected

independently from O, S and N, which heterocyclic group may bear 1 or 2 substituents selected from

oxo, hydroxy, halogeno, $C_{1-4}alkyl$, $C_{1-4}hydroxyalkyl$, $C_{1-4}alkoxy$,

$C_{1-4}alkoxyC_{1-4}alkyl$ and $C_{1-4}alkylsulphonyl$ $C_{1-4}alkyl$)), or

$(CH_2)_aY^6(CH_2)_bR^{53}$ (wherein R^{53} is as defined herein, a is 0, or an integer 1-4, b is 0 or an integer 1-4 and Y^6 represents a direct bond, $-O-$, $-C(O)-$, $-NR^{55}-$, $-NR^{50}C(O)-$ or $-C(O)NR^{57}-$ (wherein R^{55} , R^{56} , and R^{57} , which may be the same or

different, each represents hydrogen, $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$), and wherein one or more of the $(CH_2)_a$ or

$(CH_2)_b$ groups may bear one or more substituents selected from hydroxy, amino and halogeno));

with the proviso that R^5 is not alkoxy, substituted alkoxy (wherein R^5 is Y^4R^{35} and Y^4 is $-O-$ and R^{35} is $C_{1-7}alkyl$ bearing one or more substituents selected from the list given herein), $-O-$ $C_{1-7}alkanoyl$ or benzyloxy.